



Searches for User lwells (Count = 3529)

Queries 3480 through 3529.

S #	Updt	Database	Query	Time	Comment
S3529	U	USPT	6416760.pn.	2002-11-08 07:54:12	
S3528	U	USPT	irritant side effect	2002-11-07 16:13:54	
S3527	U	USPT,JPAB,EPAB,DWPI	((424/401)!.CCLS.) and (psoriasis same (alpha keto acid or pyruvic or keto octanoic or beta keto acid or acetoacetic or retinoid or retinoic or retinol or anthralin or anthranoid or dioxyanthranol or peroxide or benzoyl peroxide or minoxidil or lithium salt or antimetabolite or vitamin d or depigmentation or depigmenting or hydroquinone))	2002-11-07 12:50:50	
S3526					

U

USPT,JPAB,EPAB,DWPI

psoriasis same (alpha keto acid or pyruvic or keto octanoic or beta keto acid or acetoacetic or
retinoid or retinoic or retinol or anthralin or anthranoid or
dioxyanthranol or peroxide or benzoyl
peroxide or minoxidil or lithium salt or antimetabolite or vitamin d or
depigmentation or
depigmenting or hydroquinone)

2002-11-07

12:50:36

S3525

U

USPT,JPAB,EPAB,DWPI

(nsaid or antiinflammatory or anti inflammatory) near (alpha keto acid or pyruvic or keto
octanoic or beta keto acid or acetoacetic or retinoid or retinoic or
retinol or anthralin or
anthranoid or dioxyanthranol or peroxide or benzoyl peroxide or
minoxidil or lithium salt or
antimetabolite or vitamin d or depigmentation or depigmenting or
hydroquinone)

2002-11-07

12:38:12

S3524

U

USPT,JPAB,EPAB,DWPI

((424/401)!.CCLS.) and ((nsaid or antiinflammatory or anti inflammatory) same (alpha keto acid or pyruvic or keto octanoic or beta keto acid or acetoacetic or
retinoid or retinoic or retinol
or anthralin or anthranoid or dioxyanthranol or peroxide or benzoyl
peroxide or minoxidil or
lithium salt or antimetabolite or vitamin d or depigmentation or
depigmenting or hydroquinone
))

2002-11-07

12:37:54

S3523

U

USPT,JPAB,EPAB,DWPI

(nsaid or antiinflammatory or anti inflammatory) same (alpha keto acid or pyruvic or keto
octanoic or beta keto acid or acetoacetic or retinoid or retinoic or
retinol or anthralin or

anthranoid or dioxyanthranol or peroxide or benzoyl peroxide or
minoxidil or lithium salt or
antimetabolite or vitamin d or depigmentation or depigmenting or
hydroquinone)

2002-11-07

12:28:44

S3522

U

USPT,JPAB,EPAB,DWPI

sodium cromoglycate same (alpha keto acid or pyruvic or keto octanoic
or beta keto acid or

acetoacetic or retinoid or retinoic or retinol or anthralin or anthranoid
or dioxyanthranol or

peroxide or benzoyl peroxide or minoxidil or lithium salt or
antimetabolite or vitamin d or

depigmentation or depigmenting or hydroquinone)

2002-11-07

12:26:40

S3521

U

USPT,JPAB,EPAB,DWPI

hair and (cosmetic and ((tnf near alpha near antagonist)not (tnf near
alpha near antagonist)))

2002-11-07

12:21:59

S3520

U

USPT,JPAB,EPAB,DWPI

cosmetic and (((424/401)!.CCLS.) and ((tnf near alpha near antagonist
)not (tnf near alpha near
antagonist)))

2002-11-07

12:21:41

S3519

U

USPT,JPAB,EPAB,DWPI

((424/401)!.CCLS.) and (sodium cromoglycate)

2002-11-07

12:21:26

S3518

U

USPT,JPAB,EPAB,DWPI

sodium cromoglycate

2002-11-07

12:21:15

S3517

U

USPT,JPAB,EPAB,DWPI

((((lisophyline or a802715 or sulphasalazine)and (pharmaceutical or cosmetic))not (((tnf near alpha near antagonist)not (tnf near alpha near antagonist))or ((interleukin 1 near antagonist)and (interleukin 1 near antagonist))or ((interleukin 1 near antagonist)same (interleukin 1 near antagonist))or (cosmetic and (alpha keto acid or pyruvic or keto octanoic or beta keto acid or acetoacetic or retinoid or retinoic or retinol or anthralin or anthranoid or dioxyanthranol or peroxide or benzoyl peroxide or minoxidil or lithium salt or antimetabolite or vitamin d or depigmentation or depigmenting or hydroquinone)))) not (((auranofin or skf-105809 or lactoferin)or (lisophyline or a802715 or sulphasalazine)or (interleukin 1 near antagonist)or (tnf near alpha near antagonist)) and (alpha keto acid or pyruvic or keto octanoic or beta keto acid or acetoacetic or retinoid or retinoic or retinol or anthralin or anthranoid or dioxyanthranol or peroxide or benzoyl peroxide or minoxidil or lithium salt or antimetabolite or vitamin d or depigmentation or depigmenting or hydroquinone))

2002-11-07

11:32:47

S3516

U

USPT,JPAB,EPAB,DWPI

((((lisophyline or a802715 or sulphasalazine)and (pharmaceutical or cosmetic)) not ((((tnf near alpha near antagonist)not (tnf near alpha near antagonist)) or ((interleukin 1 near antagonist)and (interleukin 1 near antagonist)) or ((interleukin 1 near antagonist)same (interleukin 1 near antagonist)) or (cosmetic and (alpha keto acid or pyruvic or keto octanoic or beta keto acid or acetoacetic or retinoid or retinoic or retinol or anthralin or anthranoid or dioxyanthranol or peroxide or benzoyl peroxide or minoxidil or lithium salt or antimetabolite or vitamin d or depigmentation or depigmenting or hydroquinone))))

2002-11-07

11:28:01

S3515

U

USPT,JPAB,EPAB,DWPI

((lisophyline or a802715 or sulphasalazine)or (alpha keto acid or pyruvic or keto octanoic or beta keto acid or acetoacetic or retinoid or retinoic or retinol or anthralin or anthranoid or dioxyanthranol or peroxide or benzoyl peroxide or minoxidil or lithium salt or antimetabolite or vitamin d or depigmentation or depigmenting or hydroquinone)) and (pharmaceutical or cosmetic)

2002-11-07

11:27:22

S3514

U

USPT,JPAB,EPAB,DWPI

(lisophyline or a802715 or sulphasalazine) or (tnf near alpha near antagonist)

2002-11-07

11:26:10

S3513

U

USPT,JPAB,EPAB,DWPI

(auranofin or skf-105809 or lactoferin) and (lisophyline or a802715 or sulphasalazine)

2002-11-07

11:02:10

S3512

U

USPT,JPAB,EPAB,DWPI

((interleukin 1 near antagonist)and (interleukin 1 near antagonist)) not ((interleukin 1 near antagonist)same (interleukin 1 near antagonist))

2002-11-07

11:00:54

S3511

U

USPT,JPAB,EPAB,DWPI

(interleukin 1 near antagonist) and (tnf near alpha near antagonist)

2002-11-07

11:00:46

S3510

U

USPT,JPAB,EPAB,DWPI

(interleukin 1 near antagonist) same (tnf near alpha near antagonist)

2002-11-07

10:58:46

S3509

U

USPT,JPAB,EPAB,DWPI

cosmetic and (((auranofin or skf-105809 or lactoferin)or (lisophyline
or a802715 or
sulphasalazine)or (interleukin 1 near antagonist)or (tnf near alpha near
antagonist)) and (alpha
keto acid or pyruvic or keto octanoic or beta keto acid or acetoacetic or
retinoid or retinoic or
retinol or anthralin or anthranoid or dioxyanthranol or peroxide or
benzoyl peroxide or minoxidil
or lithium salt or antimetabolite or vitamin d or depigmentation or
depigmenting or
hydroquinone))

2002-11-07

10:22:37

S3508

U

USPT,JPAB,EPAB,DWPI

((auranofin or skf-105809 or lactoferin) or (lisophyline or a802715 or
sulphasalazine) or
(interleukin 1 near antagonist) or (tnf near alpha near antagonist)) and
(alpha keto acid or
pyruvic or keto octanoic or beta keto acid or acetoacetic or retinoid or
retinoic or retinol or
anthralin or anthranoid or dioxyanthranol or peroxide or benzoyl
peroxide or minoxidil or
lithium salt or antimetabolite or vitamin d or depigmentation or
depigmenting or hydroquinone)

2002-11-07

10:22:24

S3507

U

USPT,JPAB,EPAB,DWPI

tnf near alpha near antagonist

2002-11-07

10:21:57

S3506

U

USPT,JPAB,EPAB,DWPI

interleukin 1 near antagonist

2002-11-07

10:21:23

S3505

U

USPT,JPAB,EPAB,DWPI	
alpha keto acid or pyruvic or keto octanoic or beta keto acid or	
acetoacetic or retinoid or retinoic	
or retinol or anthralin or anthranoid or dioxyanthranol or peroxide or	
benzoyl peroxide or	
minoxidil or lithium salt or antimetabolite or vitamin d or	
depigmentation or depigmenting or	
hydroquinone	
	2002-11-07
	10:19:41
S3504	
U	
USPT,JPAB,EPAB,DWPI	
lisophylline or a802715 or sulphasalazine	
	2002-11-07
	10:15:18
S3503	
U	
USPT,JPAB,EPAB,DWPI	
auranofin or skf-105809 or lactoferin	
	2002-11-07
	10:14:53
S3502	
U	
USPT,JPAB,EPAB,DWPI	
benzothiophene near carboxamide	
	2002-11-07
	10:14:21
S3501	
U	
USPT,JPAB,EPAB,DWPI	
furan or benzofuran	
	2002-11-07
	10:13:52
S3500	
U	
USPT,JPAB,EPAB,DWPI	
(loratidine or cetirizine or auranofin or lisophylline or sulphasalazine or	
setastine or crotamiton	
or dexchlorpheniramine) or cinnarizine or cyclizine or	
dexchlorpheniramine or triprolidine or	
phenothiazine or promethazine or alimemazine or ebastine	
	2002-11-07
	10:13:29
S3499	
U	

USPT,JPAB,EPAB,DWPI
loratidine or cetirizine or auranofin or lisophylline or sulphasalazine or
setastine or crotamiton or
dexchlorpheniramine

2002-11-07
10:12:31

S3498

U

USPT

alpha keto acid and beta keto acid

2002-11-07
09:59:50

Welcome to STN International! Enter x:x

LOGINID:ssspta16191xw

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

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NEWS 2 Apr 08 "Ask CAS" for self-help around the clock
NEWS 3 Apr 09 BEILSTEIN: Reload and Implementation of a New Subject Area
NEWS 4 Apr 09 ZDB will be removed from STN
NEWS 5 Apr 19 US Patent Applications available in IFICDB, IFIPAT, and
IFIUDB
NEWS 6 Apr 22 Records from IP.com available in CAPLUS, HCAPLUS, and
ZCAPLUS
NEWS 7 Apr 22 BIOSIS Gene Names now available in TOXCENTER
NEWS 8 Apr 22 Federal Research in Progress (FEDRIP) now available
NEWS 9 Jun 03 New e-mail delivery for search results now available
NEWS 10 Jun 10 MEDLINE Reload
NEWS 11 Jun 10 PCTFULL has been reloaded
NEWS 12 Jul 02 FOREGE no longer contains STANDARDS file segment
NEWS 13 Jul 22 USAN to be reloaded July 28, 2002;
saved answer sets no longer valid
NEWS 14 Jul 29 Enhanced polymer searching in REGISTRY
NEWS 15 Jul 30 NETFIRST to be removed from STN
NEWS 16 Aug 08 CANCERLIT reload
NEWS 17 Aug 08 PHARMAMarketLetter(PHARMAML) - new on STN
NEWS 18 Aug 08 NTIS has been reloaded and enhanced
NEWS 19 Aug 19 Aquatic Toxicity Information Retrieval (AQUIRE)
now available on STN
NEWS 20 Aug 19 IFIPAT, IFICDB, and IFIUDB have been reloaded
NEWS 21 Aug 19 The MEDLINE file segment of TOXCENTER has been reloaded
NEWS 22 Aug 26 Sequence searching in REGISTRY enhanced
NEWS 23 Sep 03 JAPIO has been reloaded and enhanced
NEWS 24 Sep 16 Experimental properties added to the REGISTRY file
NEWS 25 Sep 16 Indexing added to some pre-1967 records in CA/CAPLUS
NEWS 26 Sep 16 CA Section Thesaurus available in CAPLUS and CA
NEWS 27 Oct 01 CASREACT Enriched with Reactions from 1907 to 1985
NEWS 28 Oct 21 EVENTLINE has been reloaded
NEWS 29 Oct 24 BEILSTEIN adds new search fields
NEWS 30 Oct 24 Nutraceuticals International (NUTRACEUT) now available on
STN
NEWS 31 Oct 25 MEDLINE SDI run of October 8, 2002

NEWS EXPRESS October 14 CURRENT WINDOWS VERSION IS V6.01,
CURRENT MACINTOSH VERSION IS V6.0a(ENG) AND V6.0Ja(JP),
AND CURRENT DISCOVER FILE IS DATED 01 OCTOBER 2002
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NEWS INTER General Internet Information
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NEWS PHONE Direct Dial and Telecommunication Network Access to STN
NEWS WWW CAS World Wide Web Site (general information)

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FILE 'HOME' ENTERED AT 09:31:24 ON 07 NOV 2002

=> fil reg

COST IN U.S. DOLLARS

SINCE FILE
ENTRY

TOTAL
SESSION

FULL ESTIMATED COST

0.21

0.21

FILE 'REGISTRY' ENTERED AT 09:31:30 ON 07 NOV 2002

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STRUCTURE FILE UPDATES: 6 NOV 2002 HIGHEST RN 471238-76-3

DICTIONARY FILE UPDATES: 6 NOV 2002 HIGHEST RN 471238-76-3

TSCA INFORMATION NOW CURRENT THROUGH MAY 20, 2002

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. See HELP PROPERTIES for more information. See STNote 27, Searching Properties in the CAS Registry File, for complete details:

<http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf>

=> e loratidine/cn

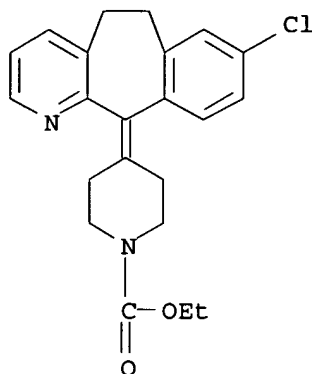
E1	1	LORATADINE/CN
E2	1	LORATADINE-PSEUDOEPHEDRINE SULFATE MIXT./CN
E3	1 -->	LORATIDINE/CN
E4	1	LORATYNE/CN
E5	1	LORATYNE D/CN
E6	1	LORAX/CN
E7	1	LORAXANTHAL MONOEPOXIDE/CN
E8	1	LORAXANTHIN DIMETHYL ETHER/CN
E9	1	LORAXANTHIN MONOEPOXIDE TRIACETATE/CN
E10	1	LORAZENE/CN
E11	1	LORAZEP/CN
E12	1	LORAZEPAM/CN

=> s e3

L1 1 LORATIDINE/CN

=> d

L1 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2002 ACS
RN 79794-75-5 REGISTRY
CN 1-Piperidinecarboxylic acid, 4-(8-chloro-5,6-dihydro-11H-benzo[5,6]cyclohepta[1,2-b]pyridin-11-ylidene)-, ethyl ester (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN 11H-Benzo[5,6]cyclohepta[1,2-b]pyridine, 1-piperidinecarboxylic acid deriv.
OTHER NAMES:
CN Anhissen
CN Bonalerg
CN Civeran
CN Claratyne
CN Claritin
CN Claritine
CN Clarityn
CN Clarityne
CN Cronopen
CN Fristamin
CN Histaloran
CN Klaritin
CN Lertamine
CN Lisino
CN Loracert
CN Loradex
CN Loranox
CN Lorastine
CN Loratadine
CN **Loratidine**
CN Loratyne
CN Lorfast
CN Lowadina
CN Optimin
CN Polaratyne
CN Pylor
CN Restamine
CN Sch 29851
CN Sensibit
CN Sohotin
CN Tadine
CN Velodan
CN Zeos
MF C22 H23 Cl N2 O2
CI COM
LC STN Files: ADISINSIGHT, ADISNEWS, AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS, BIOTECHNO, CA, CANCERLIT, CAPLUS, CASREACT, CBNB, CEN, CHEMCATS, CIN, CSCHEM, DDFU, DIOGENES, DRUGNL, DRUGPAT, DRUGU, DRUGUPDATES, EMBASE, HSDB*, IPA, MEDLINE, MRCK*, NIOSHTIC, PHAR, PHARMASEARCH, PROMT, SPECINFO, SYNTHLINE, TOXCENTER, USAN, USPAT2, USPATFULL, VETU
(*File contains numerically searchable property data)
Other Sources: WHO



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

453 REFERENCES IN FILE CA (1962 TO DATE)
 12 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
 459 REFERENCES IN FILE CAPLUS (1962 TO DATE)

=> e cetrizine/cn

E1	1	CETRIPS/CN
E2	1	CETRIZET/CN
E3	0 -->	CETRIZINE/CN
E4	1	CETRORELIX/CN
E5	1	CETRORELIX ACETATE/CN
E6	1	CETROTIDE/CN
E7	1	CETUXIMAB/CN
E8	1	CETWIST TRANSCRIPTION FACTOR (CAENORHABDITIS ELEGANS GENE
HL		H-8)/CN
E9	1	CETYL .BETA.-AMINOCROTONATE/CN
E10	1	CETYL .GAMMA.-AMINO BUTYRATE/CN
E11	1	CETYL 1,3-DIMETHYLBUTYL ETHER/CN
E12	1	CETYL 1-NAPHTHYL ETHER/CN

=> e cetirizine/cn

E1	1	CETIPRIN NOVUM/CN
E2	1	CETIR/CN
E3	1 -->	CETIRIZINE/CN
E4	1	CETIRIZINE DIHYDROCHLORIDE/CN
E5	1	CETIRIZINE HYDROCHLORIDE/CN
E6	1	CETL/CN
E7	1	CETO/CN
E8	1	CETO-2-RHODOVIBRINE/CN
E9	1	CETOBEMIDON/CN
E10	1	CETOBEMIDONE/CN
E11	1	CETOCIRE/CN
E12	1	CETOCYCLINE/CN

=> s e3

L2	1	CETIRIZINE/CN
----	---	---------------

=> d

L2 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2002 ACS
 RN 83881-51-0 REGISTRY

CN Acetic acid, [2-[4-[(4-chlorophenyl)phenylmethyl]-1-piperazinyl]ethoxy] -
(9CI) (CA INDEX NAME)

OTHER NAMES:

CN Cetirizine

FS 3D CONCORD

DR 130018-86-9

MF C21 H25 Cl N2 O3

CI COM

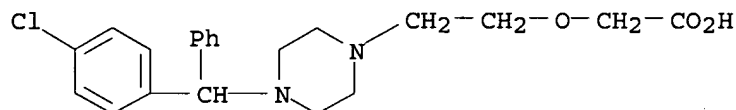
LC STN Files: ADISINSIGHT, ADISNEWS, AGRICOLA, ANABSTR, BIOBUSINESS,
BIOSIS, BIOTECHNO, CA, CANCERLIT, CAPLUS, CASREACT, CBNB, CEN,

CHEMCATS,

CIN, CSCHEM, DDFU, DIOGENES, DRUGNL, DRUGPAT, DRUGU, DRUGUPDATES,
EMBASE, IPA, MEDLINE, MRCK*, NIOSHTIC, PHAR, PROMT, RTECS*, SYNTHLINE,
TOXCENTER, USAN, USPAT2, USPATFULL, VETU

(*File contains numerically searchable property data)

Other Sources: WHO



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

430 REFERENCES IN FILE CA (1962 TO DATE)

9 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA

435 REFERENCES IN FILE CAPLUS (1962 TO DATE)

=> e auranofin/cn

E1	1	AURANAMIDE/CN
E2	1	AURANETIN/CN
E3	1 -->	AURANOFIN/CN
E4	1	AURANTHINE/CN
E5	3	AURANTIA/CN
E6	1	AURANTIACIN/CN
E7	1	AURANTIACIN, DIACETATE/CN
E8	1	AURANTIACIN, DIBENZOATE/CN
E9	1	AURANTIAZONE/CN
E10	1	AURANTIAMARIN/CN
E11	1	AURANTIAMIDE/CN
E12	1	AURANTIAMIDE ACETATE/CN

=> s e3

L3 1 AURANOFIN/CN

=> d

L3 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2002 ACS

RN 34031-32-8 REGISTRY

CN Gold, [1-(thio-.kappa.S)-.beta.-D-glucopyranose 2,3,4,6-tetraacetato](triethylphosphine)- (9CI) (CA INDEX NAME)

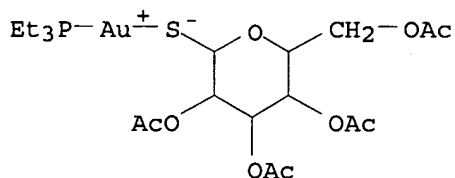
OTHER CA INDEX NAMES:

CN .beta.-D-Glucopyranose, 1-thio-, 2,3,4,6-tetraacetate, gold complex

CN Gold, (1-thio-.beta.-D-glucopyranosato)(triethylphosphine)-, 2,3,4,6-tetraacetate (8CI)

CN Gold, (1-thio-.beta.-D-glucopyranose 2,3,4,6-tetraacetato-

S) (triethylphosphine) -
 OTHER NAMES:
 CN **Auranofin**
 CN Ridaura
 CN SKF 39162
 CN SKF 39162D
 CN [(Tetra-O-acetyl-.beta.-D-glucopyranosyl)thio] (triethylphosphine)gold
 MF C20 H34 Au O9 P S
 CI CCS
 LC STN Files: ADISNEWS, AGRICOLA, ANABSTR, BIOBUSINESS, BIOSIS, BIOTECHNO,
 CA, CANCERLIT, CAPLUS, CBNB, CHEMCATS, CHEMLIST, CIN, CSCHEM, DDFU,
 DIOGENES, DRUGPAT, DRUGU, EMBASE, GMELIN*, IFICDB, IFIPAT, IFIUDB, IPA,
 MEDLINE, MRCK*, MSDS-OHS, NIOSHTIC, PHAR, PHARMASEARCH, PROMT, RTECS*,
 SYNTHLINE, TOXCENTER, USAN, USPATFULL, VETU
 (*File contains numerically searchable property data)
 Other Sources: EINECS**, WHO
 (**Enter CHEMLIST File for up-to-date regulatory information)



465 REFERENCES IN FILE CA (1962 TO DATE)
 9 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
 466 REFERENCES IN FILE CAPLUS (1962 TO DATE)

=> e lisophylline/cn
 E1 1 LISOFYLLINE/CN
 E2 1 LISOLIPIN/CN
 E3 0 --> LISOPHYLLINE/CN
 E4 1 LISOPHYLLINE/CN
 E5 1 LISORIL/CN
 E6 1 LISPAMINA/CN
 E7 1 LISPAMOL/CN
 E8 1 LISPAMOL FUMARATE/CN
 E9 1 LISPRIL/CN
 E10 1 LISSAMINE AMARANTH AC/CN
 E11 1 LISSAMINE BLUE/CN
 E12 1 LISSAMINE BLUE 2BR/CN

=> s e4
 L4 1 LISOPHYLLINE/CN

=> e
 E13 1 LISSAMINE BLUE AR/CN
 E14 1 LISSAMINE BLUE B/CN
 E15 1 LISSAMINE BLUE BF/CN
 E16 1 LISSAMINE BLUE RB/CN
 E17 1 LISSAMINE FAST RED/CN
 E18 1 LISSAMINE FAST RED 3G/CN
 E19 1 LISSAMINE FAST RED 4G/CN

E20 1 LISSAMINE FAST RED B/CN
 E21 1 LISSAMINE FAST RED BG/CN
 E22 1 LISSAMINE FAST RED BS/CN
 E23 1 LISSAMINE FAST VIOLET 2B/CN
 E24 1 LISSAMINE FAST YELLOW/CN

=> e 14

E1 4 L3T4.25/BI
 E2 4 L3T425/BI
 E3 335 --> L4/BI
 E4 1 L4,1/BI
 E5 2 L4.13.2/BI
 E6 19 L40/BI
 E7 1 L4000/BI
 E8 1 L40000/BI
 E9 1 L40001/BI
 E10 1 L40002/BI
 E11 1 L40003/BI
 E12 1 L40004/BI

=> d 14

L4 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2002 ACS

RN 100324-81-0 REGISTRY

CN 1H-Purine-2,6-dione, 3,7-dihydro-1-[(5R)-5-hydroxyhexyl]-3,7-dimethyl-
 (9CI) (CA INDEX NAME)

OTHER CA INDEX NAMES:

CN 1H-Purine-2,6-dione, 3,7-dihydro-1-(5-hydroxyhexyl)-3,7-dimethyl-, (R)-

OTHER NAMES:

CN CT 1501R

CN Lisofylline

CN **Lisophylline**

CN ProTec

FS STEREOSEARCH

MF C13 H20 N4 O3

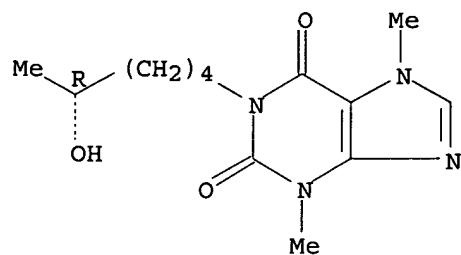
SR CA

LC STN Files: ADISINSIGHT, ADISNEWS, AGRICOLA, BEILSTEIN*, BIOBUSINESS,
 BIOSIS, BIOTECHNO, CA, CAPLUS, CASREACT, CIN, DDFU, DRUGNL, DRUGPAT,
 DRUGU, DRUGUPDATES, EMBASE, IPA, MRCK*, PHAR, PROMT, SYNTHLINE,
 TOXCENTER, USAN, USPATFULL

(*File contains numerically searchable property data)

Other Sources: WHO

Absolute stereochemistry. Rotation (-).



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

63 REFERENCES IN FILE CA (1962 TO DATE)
63 REFERENCES IN FILE CAPLUS (1962 TO DATE)

=> e sulphasalazine/cn

E1 1 SULPHARSENOLUM/CN
E2 1 SULPHARSIDE/CN
E3 1 --> SULPHASALAZINE/CN
E4 1 SULPHASIL/CN
E5 1 SULPHASOLUCIN/CN
E6 1 SULPHASOMIDINE/CN
E7 1 SULPHASOMIZOLE/CN
E8 1 SULPHAST GREEN/CN
E9 1 SULPHATASE (SALMONELLA ENTERICA TYPHI STRAIN CT18 GENE
STY24
66)/CN
E10 1 SULPHATE/CN
E11 1 SULPHATE ABC TRANSPORTER (NEISSERIA MENINGITIDIS STRAIN
C311
+ CLONE NMB0580 GENE PHO2-91 HOMOLOG)/CN
E12 1 SULPHATE ABC TRANSPORTER (NEISSERIA MENINGITIDIS STRAIN
C311
+ CLONE NMB0880 GENE PHO2-90 HOMOLOG)/CN

=> s e3

L5 1 SULPHASALAZINE/CN

=> d

L5 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2002 ACS
RN 599-79-1 REGISTRY
CN Benzoic acid, 2-hydroxy-5-[[4-[(2-pyridinylamino)sulfonyl]phenyl]azo]-
(9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Salicylic acid, 5-[[p-(2-pyridylsulfamoyl)phenyl]azo]- (6CI, 7CI, 8CI)
OTHER NAMES:
CN 2-Hydroxy-5-((4-((2-pyridinylamino)sulfonyl)phenyl)azo)benzoic acid
CN 5-[4-(2-Pyridylsulfamyl)phenylazo]-2-hydroxybenzoic acid
CN 5-[p-(2-Pyridylsulfamoyl)phenylazo]salicylic acid
CN 5-[p-(2-Pyridylsulfamyl)phenylazo]salicylic acid
CN Azopyrin
CN Azopyrine
CN Azulfidine
CN Benzosulfa
CN Reupirin
CN Salazopyridin
CN Salazopyrin
CN Salazopyrine
CN Salazosulfapyridin
CN Salazosulfapyridine
CN Salicylazosulfapyridine
CN Salisulf
CN Sulfasalazin
CN Sulfasalazine
CN **Sulphasalazine**
FS 3D CONCORD
MF C18 H14 N4 O5 S
CI COM
LC STN Files: ADISNEWS, AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS,
BIOSIS,

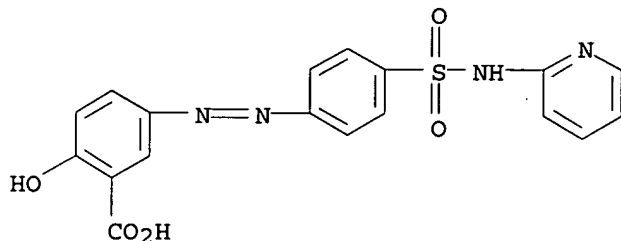
BIOTECHNO, CA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CBNB, CEN, CHEMCATS,
CHEMLIST, CIN, CSCHEM, CSNB, DDFU, DIOGENES, DRUGNL, DRUGU,
DRUGUPDATES,

EMBASE, HODOC*, HSDB*, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MRCK*,
NIOSHTIC, PHAR, PHARMASEARCH, PIRA, PROMT, RTECS*, SPECINFO, TOXCENTER,
USAN, USPATFULL, VETU

(*File contains numerically searchable property data)

Other Sources: DSL**, EINECS**, WHO

(**Enter CHEMLIST File for up-to-date regulatory information)



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

899 REFERENCES IN FILE CA (1962 TO DATE)

32 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA

903 REFERENCES IN FILE CAPLUS (1962 TO DATE)

12 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

=> e setastine/cn

E1	1	SETARON GOLDEN YELLOW RL/CN
E2	1	SETARON RED BL/CN
E3	1	--> SETASTINE/CN
E4	1	SETAVIN CAW/CN
E5	1	SETAVIN KS/CN
E6	1	SETAVIN PE/CN
E7	1	SETAVONE C/CN
E8	1	SETAVONE O/CN
E9	1	SETAVONE ST/CN
E10	1	SETAZINDOL/CN
E11	1	SETAZINE/CN
E12	1	SETCREASIN/CN

=> s e3

L6 1 SETASTINE/CN

=> d

L6 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2002 ACS

RN 64294-95-7 REGISTRY

CN 1H-Azepine, 1-[2-[1-(4-chlorophenyl)-1-phenylethoxy]ethyl]hexahydro-(9CI)

(CA INDEX NAME)

OTHER NAMES:

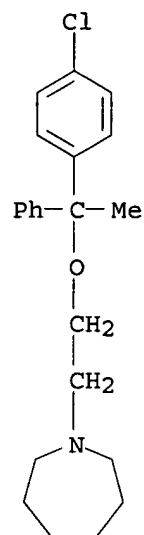
CN **Setastine**

FS 3D CONCORD

MF C22 H28 Cl N O

CI COM

LC STN Files: BEILSTEIN*, BIOSIS, CA, CAPLUS, CIN, DDFU, DRUGU, EMBASE,
 MEDLINE, MRCK*, SYNTHLINE, TOXCENTER, USAN, USPATFULL
 (*File contains numerically searchable property data)
 Other Sources: WHO



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

25 REFERENCES IN FILE CA (1962 TO DATE)
 3 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
 25 REFERENCES IN FILE CAPLUS (1962 TO DATE)

=> e crotamiton/cn

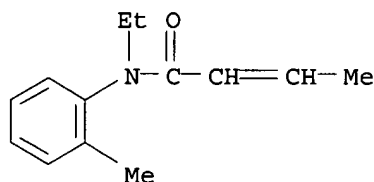
E1	1	CROTAMINE, PRO- (CROTALUS DURISSUS TERRIFICUS CLONE PCM31B
R		EDUCED) /CN
E2	1	CROTAMINE, PRO- (CROTALUS DURISSUS TERRIFICUS CLONE PCM45B
R		EDUCED) /CN
E3	1 -->	CROTAMITON/CN
E4	1	CROTAMITON-DELTAMETHRIN MIXT./CN
E5	1	CROTAMITON-PERMETHRIN MIXT./CN
E6	1	CROTAMITON-RESMETHRIN MIXT./CN
E7	1	CROTAMITON-TETRAMETHRIN MIXT./CN
E8	1	CROTAMITONE/CN
E9	1	CROTAN BK/CN
E10	1	CROTANANIC ACID/CN
E11	1	CROTANANIC ACID LACTONE/CN
E12	1	CROTANANINE/CN

=> s e3

L7 1 CROTAMITON/CN

=> d

L7 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2002 ACS
 RN 483-63-6 REGISTRY
 CN 2-Butenamide, N-ethyl-N-(2-methylphenyl)- (9CI) (CA INDEX NAME)
 OTHER CA INDEX NAMES:
 CN o-Crotonotoluidide, N-ethyl- (6CI, 7CI, 8CI)
 OTHER NAMES:
 CN Crotalgin
 CN **Crotamiton**
 CN Crotamitone
 CN Crotonyl N-ethyl-o-toluidine
 CN Eurax
 CN Euraxil
 CN N-Crotonyl-N-ethyl-o-toluidine
 CN N-Ethyl-o-crotonotoluidide
 CN Veteusan
 FS 3D CONCORD
 MF C13 H17 N O
 CI COM
 LC STN Files: ADISNEWS, AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS,
 BIOSIS,
 BIOTECHNO, CA, CABA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CBNB, CEN,
 CHEMCATS, CHEMLIST, CIN, CSCHEM, DDFU, DIOGENES, DRUGU, EMBASE, IPA,
 MEDLINE, MRCK*, PHARMASEARCH, PROMT, RTECS*, SPECINFO, TOXCENTER, USAN,
 USPATFULL, VETU
 (*File contains numerically searchable property data)
 Other Sources: EINECS**, WHO
 (**Enter CHEMLIST File for up-to-date regulatory information)



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

215 REFERENCES IN FILE CA (1962 TO DATE)
 3 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
 215 REFERENCES IN FILE CAPLUS (1962 TO DATE)
 2 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

=> e dexchlorpheniramine/cn

E1	1	DEXCEL CLEAR CONC S 380/CN
E2	1	DEXCHLORPHENIRAMINE MALEATE/CN
E3	1 -->	DEXCHLORPHENIRAMINE/CN
E4	1	DEXCHLORPHENIRAMINE HYDROGEN MALEATE/CN
E5	1	DEXCHLORPHENIRAMINE MALEATE/CN
E6	1	DEXCLAMOL/CN
E7	1	DEXCLAMOL HYDROBROMIDE/CN
E8	1	DEXCLAMOL HYDROCHLORIDE/CN
E9	1	DEXCO 2518/CN
E10	1	DEXCO 4111/CN

E11 1 DEXCO 506A/CN
E12 1 DEXCO 8508D/CN

=> s e3

L8 1 DEXCHLORPHENIRAMINE/CN

=> d

L8 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2002 ACS

RN 25523-97-1 REGISTRY

CN 2-Pyridinepropanamine, .gamma.-(4-chlorophenyl)-N,N-dimethyl-,
(.gamma.S)-

(9CI) (CA INDEX NAME)

OTHER CA INDEX NAMES:

CN 2-Pyridinepropanamine, .gamma.-(4-chlorophenyl)-N,N-dimethyl-, (S)-

CN Pyridine, 2-[p-chloro-.alpha.-[2-(dimethylamino)ethyl]benzyl]-, (S)-(+)-
(8CI)

OTHER NAMES:

CN (+)-Chlorpheniramine

CN d-Chlorpheniramine

CN **Dexchlorpheniramine**

CN S-(+)-Chlorpheniramine

CN S-Chlorpheniramine

FS STEREOSEARCH

DR 12167-38-3, 301-24-6

MF C16 H19 Cl N2

CI COM

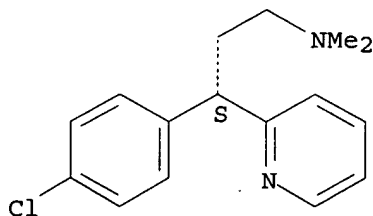
LC STN Files: ADISNEWS, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS,
BIOTECHNO, CA, CANCERLIT, CAPLUS, CASREACT, CBNB, CHEMLIST, CIN, DDFU,
DRUGU, EMBASE, HSDB*, IPA, MEDLINE, MRCK*, PHAR, PROMT, SPECINFO,
TOXCENTER, USAN, USPAT2, USPATFULL, VETU

(*File contains numerically searchable property data)

Other Sources: EINECS**, WHO

(**Enter CHEMLIST File for up-to-date regulatory information)

Absolute stereochemistry. Rotation (+).



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

217 REFERENCES IN FILE CA (1962 TO DATE)

4 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA

218 REFERENCES IN FILE CAPLUS (1962 TO DATE)

=> e benzyloxy(1)methoxy(1)tetrazol(1)benzothiophene(1)carboxamide

E1 1 BENZYLOXPHENYLACETIC/BI

E2 23201 BENZYLOXY/BI

E3 0 -->
 BENZYLOXY (L) METHOXY (L) TETRAZOL (L) BENZOTHIOPHENE (L) CARBOXAMID
 E/BI
 E4 43 BENZYLOXYACET/BI
 E5 2 BENZYLOXYACETALDEHYDE/BI
 E6 6 BENZYLOXYACETAMIDE/BI
 E7 2 BENZYLOXYACETANILIDE/BI
 E8 5 BENZYLOXYACETATE/BI
 E9 10 BENZYLOXYACETIC/BI
 E10 26 BENZYLOXYACETO/BI
 E11 2 BENZYLOXYACETOACETIC/BI
 E12 1 BENZYLOXYACETONE/BI

=> e benzyloxy(1)methoxy(1)tetrazol(1)benzothiophene(1)carboxamide/cn

E1 1 BENZYLOXY TERT-BUTYL NITROXIDE/CN
 E2 1 BENZYLOXY (ETHYL) AMINE/CN
 E3 0 -->

BENZYLOXY (L) METHOXY (L) TETRAZOL (L) BENZOTHIOPHENE (L) CARBOXAMID
 E/CN

E4 1 BENZYLOXY (PHENYLSULFONYL) METHANE/CN
 E5 1 BENZYLOXY (TERT-BUTYL) DIMETHYLSILANE/CN
 E6 1 BENZYLOXY, .ALPHA., .ALPHA.-DIMETHYL-/CN
 E7 1 BENZYLOXY, .ALPHA.-ETHYL-.ALPHA.-METHYL-/CN
 E8 1 BENZYLOXY, .ALPHA.-METHYL-/CN
 E9 1 BENZYLOXY, DIHYDROXY-/CN
 E10 1 BENZYLOXY, P-METHYL-/CN
 E11 1 BENZYLOXY-1-BROMO-2-FLUOROBENZENE/CN
 E12 1 BENZYLOXY-1-NAPHTHYLPHENYLSILANE/CN

=> log y

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

47.30

47.51

STN INTERNATIONAL LOGOFF AT 09:35:44 ON 07 NOV 2002